

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020924\
 Data File : BP019650.D
 Acq On : 10 Feb 2024 01:14
 Operator : MA/JU
 Sample : P1413-03
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 ETGI-347

Quant Time: Feb 10 03:14:20 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 01 00:51:48 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.905	152	99186	20.000	ng	-0.01
21) Naphthalene-d8	10.704	136	341906	20.000	ng	-0.01
39) Acenaphthene-d10	14.540	164	188098	20.000	ng	-0.01
64) Phenanthrene-d10	17.298	188	415846	20.000	ng	0.00
76) Chrysene-d12	21.392	240	492456	20.000	ng	0.00
86) Perylene-d12	23.810	264	574810	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.481	112	584995	95.070	ng	0.00
7) Phenol-d6	7.081	99	746690	89.997	ng	0.00
23) Nitrobenzene-d5	9.075	82	515623	65.347	ng	-0.01
42) 2,4,6-Tribromophenol	16.034	330	322461	117.414	ng	0.00
45) 2-Fluorobiphenyl	13.169	172	1026460	67.462	ng	-0.01
79) Terphenyl-d14	19.863	244	1806646	60.332	ng	0.00
Target Compounds						Qvalue
49) Acenaphthylene	14.263	152	157889	9.132	ng	98
52) Acenaphthene	14.604	154	102006	9.506	ng	98
55) Dibenzofuran	14.940	168	114710	6.563	ng	99
58) Fluorene	15.592	166	171267	12.303	ng	98
71) Phenanthrene	17.339	178	1031013	47.111	ng	100
72) Anthracene	17.433	178	743221	34.044	ng	100
73) Carbazole	17.704	167	99433	5.014	ng	98
75) Fluoranthene	19.310	202	2602929	97.363	ng	99
78) Pyrene	19.663	202	2196341	63.194	ng	99
81) Benzo(a)anthracene	21.374	228	1322431	38.149	ng	96
83) Chrysene	21.427	228	1391842	42.290	ng	98
85) Di-n-octyl phthalate	22.316	149	69833	2.008	ng	# 69
87) Indeno(1,2,3-cd)pyrene	26.333	276	352193	8.004	ng	97
88) Benzo(b)fluoranthene	23.080	252	1446496	39.928	ng	99
89) Benzo(k)fluoranthene	23.121	252	394846m	11.410	ng	
90) Benzo(a)pyrene	23.704	252	689915	21.403	ng	97
91) Dibenzo(a,h)anthracene	26.345	278	109078	3.014	ng	# 87
92) Benzo(g,h,i)perylene	27.115	276	301184	8.204	ng	98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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