

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042721\
 Data File : BM029682.D
 Acq On : 27 Apr 2021 20:16
 Operator : CG/JU
 Sample : M2176-04DL 2X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 STOCKPILE-COMP-1DL

Manual Integrations
 APPROVED

mohammad
 4/28/2021 11:00:42 AM

Quant Time: Apr 28 01:22:57 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM042121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 26 15:01:01 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.610	152	85595	20.00	ng	-0.01
21) Naphthalene-d8	10.392	136	307125	20.00	ng	0.00
39) Acenaphthene-d10	14.263	164	209403	20.00	ng	0.00
64) Phenanthrene-d10	17.004	188	477197	20.00	ng	0.00
76) Chrysene-d12	21.192	240	558005	20.00	ng	0.00
86) Perylene-d12	23.380	264	583182	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.210	112	169594	39.72	ng	0.00
7) Phenol-d6	6.793	99	209608	32.87	ng	0.00
23) Nitrobenzene-d5	8.769	82	140422	23.55	ng	0.00
42) 2,4,6-Tribromophenol	15.751	330	111314	38.44	ng	-0.01
45) 2-Fluorobiphenyl	12.875	172	388008	25.15	ng	-0.01
79) Terphenyl-d14	19.639	244	616633	20.95	ng	0.00
Target Compounds						Qvalue
49) Acenaphthylene	13.980	152	81116	4.32	ng	99
71) Phenanthrene	17.051	178	535734	19.50	ng	99
72) Anthracene	17.139	178	193709	7.15	ng	99
75) Fluoranthene	19.068	202	1266339	38.01	ng	98
78) Pyrene	19.433	202	1099907	30.64	ng	99
81) Benzo(a)anthracene	21.174	228	664482	18.00	ng	97
83) Chrysene	21.227	228	558657	15.40	ng	97
87) Indeno(1,2,3-cd)pyrene	25.568	276	384309	8.52	ng	# 95
88) Benzo(b)fluoranthene	22.727	252	810437m	21.31	ng	
89) Benzo(k)fluoranthene	22.762	252	240922m	6.32	ng	
90) Benzo(a)pyrene	23.286	252	589157	16.50	ng	98
91) Dibenzo(a,h)anthracene	25.574	278	96729	2.49	ng	# 85
92) Benzo(g,h,i)perylene	26.238	276	374942	10.43	ng	99

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

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