

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112020\
 Data File : BP004025.D
 Acq On : 20 Nov 2020 15:15
 Operator : CG/JU
 Sample : L4716-02
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 STOCKPILE-2

Manual Integrations
 APPROVED

mohammad
 11/24/2020 8:55:16 AM

Quant Time: Nov 20 16:22:49 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP110320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 20 12:19:10 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.269	152	131828	20.00	ng	0.00
21) Naphthalene-d8	11.099	136	510463	20.00	ng	# 0.00
39) Acenaphthene-d10	14.887	164	305931	20.00	ng	0.00
64) Phenanthrene-d10	17.622	188	632902	20.00	ng	0.00
76) Chrysene-d12	21.657	240	593654	20.00	ng	0.00
86) Perylene-d12	24.127	264	588822	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.805	112	759888	96.20	ng	0.00
7) Phenol-d6	7.452	99	999496	88.15	ng	0.00
23) Nitrobenzene-d5	9.428	82	656977	63.03	ng	-0.01
42) 2,4,6-Tribromophenol	16.375	330	389483	101.79	ng	0.00
45) 2-Fluorobiphenyl	13.510	172	1451973	66.35	ng	0.00
79) Terphenyl-d14	20.122	244	1860430	60.55	ng	0.00
Target Compounds						Qvalue
10) Phenol	7.475	94	84914	7.76	ng	77
50) Dimethylphthalate	14.316	163	88239	3.69	ng	98
71) Phenanthrene	17.663	178	164818	4.64	ng	98
75) Fluoranthene	19.598	202	541655	13.22	ng	100
78) Pyrene	19.945	202	509498	12.65	ng	98
81) Benzo(a)anthracene	21.639	228	276033	7.05	ng	97
83) Chrysene	21.692	228	278605	7.41	ng	97
87) Indeno(1,2,3-cd)pyrene	26.674	276	173986	4.10	ng	# 93
88) Benzo(b)fluoranthene	23.375	252	361278m	9.38	ng	
89) Benzo(k)fluoranthene	23.416	252	136404m	3.59	ng	
90) Benzo(a)pyrene	24.016	252	246959	6.92	ng	98
92) Benzo(g,h,i)perylene	27.457	276	179388	5.23	ng	96

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

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