

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031522\
 Data File : BP009418.D
 Acq On : 16 Mar 2022 02:23
 Operator : CG/JU
 Sample : N1886-04DL 5X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 30970DL

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 03/16/2022
 Supervised By :Jagrut Upadhyay 03/16/2022

Quant Time: Mar 16 03:23:38 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Mar 05 01:55:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.816	152	357042	20.000	ng	0.00
21) Naphthalene-d8	10.604	136	1363634	20.000	ng	-0.01
39) Acenaphthene-d10	14.457	164	914069	20.000	ng	0.00
64) Phenanthrene-d10	17.216	188	1911429	20.000	ng	0.00
76) Chrysene-d12	21.327	240	1930114	20.000	ng	0.00
86) Perylene-d12	23.745	264	2151555	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.422	112	161015	6.583	ng	0.01
7) Phenol-d6	7.011	99	118710	3.819	ng	0.02
23) Nitrobenzene-d5	8.969	82	350477	14.156	ng	0.00
42) 2,4,6-Tribromophenol	15.957	330	276477	25.049	ng	0.00
45) 2-Fluorobiphenyl	13.075	172	854924	12.862	ng	-0.01
79) Terphenyl-d14	19.792	244	1147014	11.215	ng	0.00
Target Compounds						Qvalue
10) Phenol	7.040	94	100469	3.121	ng	95
20) 3+4-Methylphenols	8.605	107	966991	34.630	ng	90
27) 2,4-Dimethylphenol	9.810	122	116293m	5.202	ng	
70) Pentachlorophenol	16.881	266	91236	6.267	ng	98
84) Bis(2-ethylhexyl)phtha...	21.245	149	214741	2.855	ng	# 96

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

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