

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071122\
 Data File : BP010953.D
 Acq On : 11 Jul 2022 17:34
 Operator : CG/JU
 Sample : PB146134BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK134

Quant Time: Jul 11 18:26:05 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jul 06 23:13:53 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.699	152	468115	20.000	ng/u1	0.00
20) Naphthalene-d8	10.493	136	1864065	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.357	164	910240	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.122	188	1647157	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.233	240	1608954	20.000	ng/u1	# 0.00
88) Perylene-d12	23.592	264	1806272	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.199	96	78671	6.357	ng/uL	0.00
4) Pyridine-d5	3.605	84	857304	26.026	ng/u1	0.00
7) Phenol-d5	6.875	99	1149788	31.028	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.040	67	744620	30.726	ng/u1	0.00
11) 2-Chlorophenol-d4	7.234	132	958660	31.182	ng/u1	0.00
15) 4-Methylphenol-d8	8.410	113	876178	29.910	ng/u1	-0.01
21) Nitrobenzene-d5	8.863	128	419672	35.366	ng/u1	0.00
24) 2-Nitrophenol-d4	9.581	143	386650	39.384	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.116	165	707143	29.151	ng/u1	0.00
31) 4-Chloroaniline-d4	10.640	131	910479	22.994	ng/u1	0.00
46) Dimethylphthalate-d6	13.775	166	1758960	28.210	ng/u1	0.00
49) Acenaphthylene-d8	14.045	160	2311286	29.044	ng/u1	0.00
54) 4-Nitrophenol-d4	14.569	143	285587	29.324	ng/u1	0.00
60) Fluorene-d10	15.351	176	1563455	29.219	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.475	200	151064	27.682	ng/u1	0.00
73) Anthracene-d10	17.216	188	2251564	30.724	ng/u1	0.00
81) Pyrene-d10	19.469	212	2477095	28.259	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.439	264	2805606	31.432	ng/u1	0.00

Target Compounds Qvalue

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

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