

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
 Data File : BP013268.D
 Acq On : 23 Dec 2022 08:31
 Operator : CG/JU
 Sample : PB149784BL
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK784

Quant Time: Dec 24 00:22:20 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Dec 24 00:19:17 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.951	152	514308	20.000	ng/u1	0.00
20) Naphthalene-d8	10.763	136	2058912	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.598	164	1126685	20.000	ng/u1	0.01
64) Phenanthrene-d10	17.357	188	2212862	20.000	ng/u1	0.01
79) Chrysene-d12	21.451	240	2326679	20.000	ng/u1	0.02
88) Perylene-d12	23.939	264	2712344	20.000	ng/u1	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.334	96	65801	5.470	ng/uL	0.00
4) Pyridine-d5	3.763	84	960809	28.119	ng/u1	0.00
7) Phenol-d5	7.110	99	1199310	28.629	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	744851	29.475	ng/u1	0.00
11) 2-Chlorophenol-d4	7.475	132	966937	29.321	ng/u1	0.00
15) 4-Methylphenol-d8	8.663	113	921285	26.840	ng/u1	0.00
21) Nitrobenzene-d5	9.116	128	458197	30.289	ng/u1	0.00
24) 2-Nitrophenol-d4	9.845	143	514411	30.205	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	915822	27.686	ng/u1	0.00
31) 4-Chloroaniline-d4	10.904	131	1221056	26.147	ng/u1	0.00
46) Dimethylphthalate-d6	14.004	166	2458906	28.397	ng/u1	0.00
49) Acenaphthylene-d8	14.292	160	2817736	29.614	ng/u1	0.01
54) 4-Nitrophenol-d4	14.798	143	328676	21.101	ng/u1	0.00
60) Fluorene-d10	15.592	176	2040965	27.861	ng/u1	0.01
65) 4,6-Dinitro-2-methylph...	15.716	200	230224	16.167	ng/u1	0.01
73) Anthracene-d10	17.457	188	3045106	30.502	ng/u1	0.01
81) Pyrene-d10	19.692	212	3734184	27.961	ng/u1	0.01
92) Benzo(a)pyrene-d12	23.780	264	4130652	30.553	ng/u1	0.02

Target Compounds Qvalue

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

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