

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
 Data File : BP013281.D
 Acq On : 23 Dec 2022 16:01
 Operator : CG/JU
 Sample : PB149810BL
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK810

Quant Time: Dec 24 00:25:58 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	7.940	152	552147	20.000	ng/u1	0.00
20) Naphthalene-d8	10.758	136	2273671	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.587	164	1312793	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.345	188	2297071	20.000	ng/u1	0.00
79) Chrysene-d12	21.439	240	2110964	20.000	ng/u1	0.00
88) Perylene-d12	23.927	264	2445937	20.000	ng/u1	0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.334	96	70625	5.469	ng/uL	0.00
4) Pyridine-d5	3.758	84	1021887	27.857	ng/u1	0.00
7) Phenol-d5	7.105	99	1296074	28.818	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	788329	29.058	ng/u1	0.00
11) 2-Chlorophenol-d4	7.469	132	1048521	29.616	ng/u1	0.00
15) 4-Methylphenol-d8	8.658	113	1019594	27.669	ng/u1	0.00
21) Nitrobenzene-d5	9.111	128	506339	30.310	ng/u1	0.00
24) 2-Nitrophenol-d4	9.834	143	567818	30.192	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.375	165	1039308	28.451	ng/u1	0.00
31) 4-Chloroaniline-d4	10.899	131	1386888	26.893	ng/u1	0.00
46) Dimethylphthalate-d6	13.993	166	2896897	28.712	ng/u1	0.00
49) Acenaphthylene-d8	14.281	160	3240262	29.227	ng/u1	0.00
54) 4-Nitrophenol-d4	14.793	143	340333	18.752	ng/u1	0.00
60) Fluorene-d10	15.581	176	2328110	27.275	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.704	200	250273	16.931	ng/u1	0.00
73) Anthracene-d10	17.445	188	3133652	30.238	ng/u1	0.00
81) Pyrene-d10	19.681	212	3494160	28.837	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.769	264	3728987	30.587	ng/u1	0.01

Target Compounds	Qvalue

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

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