

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122722\
 Data File : BP013337.D
 Acq On : 27 Dec 2022 14:18
 Operator : CG/JU
 Sample : PB149856BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK856

Quant Time: Dec 28 00:18:27 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 28 00:16:35 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.928	152	416869	20.000	ng/u1	0.00
20) Naphthalene-d8	10.746	136	1792540	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.575	164	1249837	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.333	188	2953525	20.000	ng/u1	0.00
79) Chrysene-d12	21.421	240	2778524	20.000	ng/u1	0.00
88) Perylene-d12	23.892	264	2621953	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.322	96	57025	5.849	ng/uL	0.00
4) Pyridine-d5	3.746	84	781597	28.220	ng/u1	0.00
7) Phenol-d5	7.093	99	984028	28.980	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.257	67	645514	31.515	ng/u1	0.00
11) 2-Chlorophenol-d4	7.457	132	798347	29.867	ng/u1	0.00
15) 4-Methylphenol-d8	8.646	113	807982	29.041	ng/u1	0.00
21) Nitrobenzene-d5	9.099	128	406793	30.887	ng/u1	0.00
24) 2-Nitrophenol-d4	9.822	143	461946	31.155	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.363	165	822825	28.571	ng/u1	0.00
31) 4-Chloroaniline-d4	10.881	131	1190364	29.278	ng/u1	0.00
46) Dimethylphthalate-d6	13.981	166	3089736	32.166	ng/u1	0.00
49) Acenaphthylene-d8	14.269	160	3161810	29.956	ng/u1	0.00
54) 4-Nitrophenol-d4	14.781	143	451451	26.128	ng/u1	0.00
60) Fluorene-d10	15.569	176	2575815	31.697	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.692	200	442733	23.294	ng/u1	0.00
73) Anthracene-d10	17.433	188	4267637	32.028	ng/u1	0.00
81) Pyrene-d10	19.669	212	5226500	32.771	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.733	264	4206178	32.185	ng/u1	0.00

Target Compounds Qvalue

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

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