

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082323\
 Data File : BP016723.D
 Acq On : 23 Aug 2023 19:49
 Operator : MA/JU
 Sample : PB154647BL
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK647

Quant Time: Aug 24 00:56:19 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081423.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Aug 23 18:42:31 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.040	152	340256	20.000	ng/u1	0.00
20) Naphthalene-d8	10.869	136	1528606	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.693	164	954732	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.457	188	2126773	20.000	ng/u1	0.00
79) Chrysene-d12	21.557	240	1629087	20.000	ng/u1	0.00
88) Perylene-d12	24.133	264	1493811	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	75352	8.415	ng/uL	0.00
4) Pyridine-d5	3.823	84	891240	32.714	ng/u1	0.00
7) Phenol-d5	7.181	99	1024562	32.662	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.375	67	666968	32.281	ng/u1	0.00
11) 2-Chlorophenol-d4	7.558	132	733857	32.571	ng/u1	0.00
15) 4-Methylphenol-d8	8.746	113	798476	32.012	ng/u1	0.00
21) Nitrobenzene-d5	9.240	128	378531	35.193	ng/u1	0.00
24) 2-Nitrophenol-d4	9.952	143	401599	39.312	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.475	165	684940	34.070	ng/u1	0.00
31) 4-Chloroaniline-d4	11.028	131	1102532	31.963	ng/u1	0.00
46) Dimethylphthalate-d6	14.104	166	2350069	34.032	ng/u1	0.00
49) Acenaphthylene-d8	14.393	160	2640693	33.880	ng/u1	0.00
54) 4-Nitrophenol-d4	14.893	143	384329	27.320	ng/u1	0.00
60) Fluorene-d10	15.687	176	2004434	34.048	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.804	200	258707	25.570	ng/u1	0.00
73) Anthracene-d10	17.557	188	3194956	34.730	ng/u1	0.00
81) Pyrene-d10	19.792	212	3498655	40.987	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.963	264	2517054	35.160	ng/u1	0.00

Target Compounds Qvalue

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

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