

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082223\
 Data File : BP016701.D
 Acq On : 22 Aug 2023 13:26
 Operator : MA/JU
 Sample : PB154937BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK937

Quant Time: Aug 23 01:15:23 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081423.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Aug 17 01:38:34 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.046	152	345723	20.000	ng/u1	0.00
20) Naphthalene-d8	10.875	136	1582193	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.692	164	1003944	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.463	188	2271989	20.000	ng/u1	0.00
79) Chrysene-d12	21.557	240	1874458	20.000	ng/u1	0.00
88) Perylene-d12	24.133	264	1778468	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	69420	7.630	ng/uL	0.00
4) Pyridine-d5	3.822	84	851382	30.757	ng/u1	0.00
7) Phenol-d5	7.181	99	973369	30.539	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.381	67	634566	30.228	ng/u1	0.00
11) 2-Chlorophenol-d4	7.563	132	695014	30.359	ng/u1	0.00
15) 4-Methylphenol-d8	8.745	113	767767	30.294	ng/u1	-0.01
21) Nitrobenzene-d5	9.240	128	369800	33.217	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.957	143	389075	36.796	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.481	165	660832	31.758	ng/u1	-0.01
31) 4-Chloroaniline-d4	11.028	131	1091435	30.570	ng/u1	-0.01
46) Dimethylphthalate-d6	14.104	166	2368682	32.620	ng/u1	0.00
49) Acenaphthylene-d8	14.392	160	2630814	32.099	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.898	143	415235	28.070	ng/u1	0.00
60) Fluorene-d10	15.692	176	2032905	32.839	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.810	200	274443	25.391	ng/u1	0.00
73) Anthracene-d10	17.563	188	3250590	33.076	ng/u1	0.00
81) Pyrene-d10	19.792	212	3680961	37.478	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.968	264	2839376	33.314	ng/u1	0.00

Target Compounds	Qvalue

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

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