

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082123\
 Data File : BP016682.D
 Acq On : 21 Aug 2023 19:39
 Operator : MA/JU
 Sample : PB154953BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK953

Quant Time: Aug 22 00:43:45 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	306316	20.000	ng/u1	0.00
20) Naphthalene-d8	10.875	136	1401862	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.698	164	884415	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.463	188	1955544	20.000	ng/u1	0.00
79) Chrysene-d12	21.557	240	1539216	20.000	ng/u1	0.00
88) Perylene-d12	24.139	264	1390674	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	56351	6.990	ng/uL	0.00
4) Pyridine-d5	3.822	84	877129	35.764	ng/u1	0.00
7) Phenol-d5	7.187	99	1049549	37.165	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.381	67	665139	35.760	ng/u1	0.00
11) 2-Chlorophenol-d4	7.563	132	744461	36.703	ng/u1	0.00
15) 4-Methylphenol-d8	8.752	113	825487	36.762	ng/u1	0.00
21) Nitrobenzene-d5	9.246	128	387724	39.307	ng/u1	0.00
24) 2-Nitrophenol-d4	9.957	143	407728	43.521	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.481	165	725207	39.335	ng/u1	-0.01
31) 4-Chloroaniline-d4	11.034	131	1158128	36.611	ng/u1	0.00
46) Dimethylphthalate-d6	14.110	166	2452701	38.342	ng/u1	0.00
49) Acenaphthylene-d8	14.398	160	2770570	38.372	ng/u1	0.00
54) 4-Nitrophenol-d4	14.898	143	419636	32.202	ng/u1	0.00
60) Fluorene-d10	15.692	176	2100306	38.513	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.810	200	269434	28.962	ng/u1	0.00
73) Anthracene-d10	17.563	188	3460368	40.909	ng/u1	0.00
81) Pyrene-d10	19.792	212	3845225	47.677	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.968	264	2761795	41.440	ng/u1	0.00

Target Compounds Qvalue

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

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