

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101723\
 Data File : BP017701.D
 Acq On : 17 Oct 2023 16:55
 Operator : MA/JU
 Sample : PB156336BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK336

Quant Time: Oct 18 00:58:37 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 12 16:14:36 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.899	152	88355	20.000	ng/u1	0.00
20) Naphthalene-d8	10.734	136	363554	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.599	164	229510	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.392	188	528824	20.000	ng/u1	0.00
79) Chrysene-d12	21.533	240	579739	20.000	ng/u1	-0.01
88) Perylene-d12	24.104	264	736169	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	14859	6.200	ng/uL	0.00
4) Pyridine-d5	3.693	84	201588	30.430	ng/u1	0.00
7) Phenol-d5	7.058	99	245442	29.838	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.246	67	161001	32.274	ng/u1	0.00
11) 2-Chlorophenol-d4	7.417	132	206102	32.369	ng/u1	0.00
15) 4-Methylphenol-d8	8.622	113	195428	29.744	ng/u1	0.00
21) Nitrobenzene-d5	9.111	128	94058	30.802	ng/u1	0.00
24) 2-Nitrophenol-d4	9.828	143	102131	29.741	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.358	165	181720	28.156	ng/u1	0.00
31) 4-Chloroaniline-d4	10.910	131	275354	29.764	ng/u1	0.00
46) Dimethylphthalate-d6	14.016	166	633499	31.870	ng/u1	0.00
49) Acenaphthylene-d8	14.299	160	668478	31.230	ng/u1	0.00
54) 4-Nitrophenol-d4	14.857	143	77590	23.340	ng/u1	0.00
60) Fluorene-d10	15.604	176	564193	33.368	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.751	200	80466	21.528	ng/u1	0.00
73) Anthracene-d10	17.492	188	929586	34.051	ng/u1	0.00
81) Pyrene-d10	19.751	212	1182187	33.881	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.933	264	1383309	33.472	ng/u1	0.00

Target Compounds	Qvalue

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

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