

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040724\
 Data File : BM045006.D
 Acq On : 07 Apr 2024 14:19
 Operator : MA/JU
 Sample : PB160068BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK068

Quant Time: Apr 08 03:47:19 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Apr 08 01:45:36 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.628	152	77196	20.000	ng/u1	0.00
20) Naphthalene-d8	10.404	136	330452	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.275	164	204069	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.039	188	427945	20.000	ng/u1	0.00
79) Chrysene-d12	21.245	240	382982	20.000	ng/u1	0.00
88) Perylene-d12	23.462	264	459241	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	12606	6.098	ng/uL	0.00
4) Pyridine-d5	3.605	84	165699	29.475	ng/u1	0.00
7) Phenol-d5	6.810	99	205681	31.429	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.981	67	125128	32.092	ng/u1	0.00
11) 2-Chlorophenol-d4	7.163	132	171275	33.406	ng/u1	0.00
15) 4-Methylphenol-d8	8.340	113	169724	33.133	ng/u1	0.00
21) Nitrobenzene-d5	8.793	128	83697	32.974	ng/u1	0.00
24) 2-Nitrophenol-d4	9.504	143	90510	33.783	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.034	165	160215	32.616	ng/u1	0.00
31) 4-Chloroaniline-d4	10.563	131	273661	34.992	ng/u1	0.00
46) Dimethylphthalate-d6	13.704	166	483737	34.045	ng/u1	0.00
49) Acenaphthylene-d8	13.969	160	585368	34.873	ng/u1	0.00
54) 4-Nitrophenol-d4	14.516	143	74506	25.125	ng/u1	0.00
60) Fluorene-d10	15.280	176	439040	34.659	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.421	200	54996	21.674	ng/u1	0.00
73) Anthracene-d10	17.133	188	690828	35.885	ng/u1	-0.01
81) Pyrene-d10	19.439	212	778827	41.300	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.321	264	796690	35.562	ng/u1	-0.01

Target Compounds Qvalue

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

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