

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072722\
 Data File : BM035965.D
 Acq On : 28 Jul 2022 00:30
 Operator : CG/JU
 Sample : PB146419BL
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK419

Quant Time: Jul 28 01:03:07 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM072522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jul 27 23:55:50 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.881	152	345581	20.000	ng/u1	0.00
20) Naphthalene-d8	10.687	136	1465299	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.522	164	876092	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.257	188	1706012	20.000	ng/u1	0.00
79) Chrysene-d12	21.433	240	1495919	20.000	ng/u1	0.00
88) Perylene-d12	23.809	264	1500792	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	42691	4.649	ng/uL	0.00
4) Pyridine-d5	3.693	84	570802	22.687	ng/u1	0.00
7) Phenol-d5	7.046	99	702936	24.256	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.216	67	452144	23.404	ng/u1	0.00
11) 2-Chlorophenol-d4	7.410	132	624173	25.948	ng/u1	0.00
15) 4-Methylphenol-d8	8.598	113	565584	24.091	ng/u1	0.00
21) Nitrobenzene-d5	9.051	128	304402	26.216	ng/u1	0.00
24) 2-Nitrophenol-d4	9.769	143	319939	28.282	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.310	165	524017	25.612	ng/u1	0.00
31) 4-Chloroaniline-d4	10.828	131	826635	23.866	ng/u1	0.00
46) Dimethylphthalate-d6	13.933	166	1618915	27.128	ng/u1	0.00
49) Acenaphthylene-d8	14.216	160	2050056	26.924	ng/u1	0.00
54) 4-Nitrophenol-d4	14.722	143	269437	21.986	ng/u1	0.00
60) Fluorene-d10	15.510	176	1399153	25.945	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.633	200	203144	22.152	ng/u1	0.00
73) Anthracene-d10	17.357	188	2097860	27.199	ng/u1	0.00
81) Pyrene-d10	19.645	212	2318332	27.963	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.650	264	2105579	27.635	ng/u1	0.00

Target Compounds Qvalue

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

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