

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080924\
 Data File : BM047124.D
 Acq On : 09 Aug 2024 17:01
 Operator : RC/JU
 Sample : PB162595BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK595

Quant Time: Aug 10 01:16:55 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Aug 09 01:30:34 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.386	152	189356	20.000	ng/u1	0.00
20) Naphthalene-d8	10.139	136	749308	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.039	164	463677	20.000	ng/u1	0.00
64) Phenanthrene-d10	16.798	188	935774	20.000	ng/u1	0.00
79) Chrysene-d12	21.033	240	775710	20.000	ng/u1	# 0.00
88) Perylene-d12	23.732	264	877030	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.016	96	30905	6.466	ng/uL	0.00
4) Pyridine-d5	3.404	84	439014	33.674	ng/u1	0.00
7) Phenol-d5	6.592	99	518049	36.015	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.745	67	335468	35.444	ng/u1	0.00
11) 2-Chlorophenol-d4	6.928	132	406525	33.363	ng/u1	0.00
15) 4-Methylphenol-d8	8.104	113	402728	35.342	ng/u1	0.00
21) Nitrobenzene-d5	8.528	128	200510	33.347	ng/u1	0.00
24) 2-Nitrophenol-d4	9.239	143	220654	33.604	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	9.780	165	397048	31.495	ng/u1	0.00
31) 4-Chloroaniline-d4	10.292	131	556923	33.613	ng/u1	0.00
46) Dimethylphthalate-d6	13.468	166	1140396	32.166	ng/u1	0.00
49) Acenaphthylene-d8	13.727	160	1344771	33.981	ng/u1	0.00
54) 4-Nitrophenol-d4	14.286	143	195397	28.505	ng/u1	0.00
60) Fluorene-d10	15.045	176	980632	32.334	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.186	200	145180	24.556	ng/u1	0.00
73) Anthracene-d10	16.898	188	1504743	33.920	ng/u1	0.00
81) Pyrene-d10	19.215	212	1631637	36.744	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.556	264	1585497	34.332	ng/u1	0.00

Target Compounds Qvalue

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

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