

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082123\
 Data File : BP016683.D
 Acq On : 21 Aug 2023 20:13
 Operator : MA/JU
 Sample : PB154937BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK937

Quant Time: Aug 22 00:45:51 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	338704	20.000	ng/u1	0.00
20) Naphthalene-d8	10.875	136	1568474	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.698	164	978790	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.463	188	2160036	20.000	ng/u1	0.00
79) Chrysene-d12	21.557	240	1738381	20.000	ng/u1	0.00
88) Perylene-d12	24.133	264	1570168	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	97684	10.959	ng/uL	0.00
4) Pyridine-d5	3.823	84	1242150	45.804	ng/u1	0.00
7) Phenol-d5	7.187	99	1406832	45.053	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.381	67	901497	43.833	ng/u1	0.00
11) 2-Chlorophenol-d4	7.563	132	994512	44.342	ng/u1	0.00
15) 4-Methylphenol-d8	8.746	113	1114824	44.900	ng/u1	-0.01
21) Nitrobenzene-d5	9.246	128	529297	47.960	ng/u1	0.00
24) 2-Nitrophenol-d4	9.957	143	564582	53.862	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.481	165	974423	47.238	ng/u1	-0.01
31) 4-Chloroaniline-d4	11.034	131	1577663	44.575	ng/u1	0.00
46) Dimethylphthalate-d6	14.110	166	3280569	46.339	ng/u1	0.00
49) Acenaphthylene-d8	14.398	160	3714731	46.488	ng/u1	0.00
54) 4-Nitrophenol-d4	14.898	143	592057	41.052	ng/u1	0.00
60) Fluorene-d10	15.692	176	2811590	46.585	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.810	200	387332	37.693	ng/u1	0.00
73) Anthracene-d10	17.563	188	4433830	47.455	ng/u1	0.00
81) Pyrene-d10	19.798	212	4904214	53.841	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.968	264	3649996	48.506	ng/u1	0.00

Target Compounds Qvalue

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

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